

On the quantum discord of general X states

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Abstract Quantum discord Q is a function of density matrix elements. The domain of such a function in the case of two-qubit system with X density matrix may consist of three subdomains at most: two ones where the quantum discord is expressed in closed analytical forms ($Q_{\pi/2}$ and Q_0) and an intermediate subdomain for which, to extract the quantum discord Q_θ , it is required to solve in general numerically a one-dimensional minimization problem to find the optimal measurement angle $\theta \in (0, \pi/2)$. Hence the quantum discord is given by a piecewise-analytic-numerical formula $Q = \min\{Q_{\pi/2}, Q_\theta, Q_0\}$. Equations for determining the boundaries between these subdomains are obtained. The boundaries consist of bifurcation points. The Q_θ subdomains are discovered in the generalized Horodecki states, in the dynamical phase flip channel model, in the anisotropic spin systems at thermal equilibrium, in the heteronuclear dimers in an external magnetic field. We found that transitions between Q_θ subdomain and $Q_{\pi/2}$ and Q_0 ones occur suddenly but continuously and smoothly, i.e., nonanalyticity is hidden and can be observed in higher derivatives of discord function.

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1 Introduction

At present we have a situation where further miniaturization of electronics will inevitably lead to molecular size components. Designing such components

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requires application of the laws of quantum mechanics. This is expected to lead to a technological breakthrough which will be achieved through employing the holy of holies of the quantum theory — so-called quantum correlations.

Initially, the entanglement has been considered as a quantum correlation [1,2]. Quantum entanglement is able to bind different parts of systems, even in the case when there is no interaction between those parts (the Einstein-Podolsky-Rosen effect). Quantum entanglement exists only in nonseparable states of bi- and multipartite systems. However, it appears in the last years that there are quantum correlations more general and more fundamental than entanglement. In particular, they can be present in certain separable states, i.e., when the quantum entanglement is absent. As a measure of total purely quantum correlations in bipartite systems, the quantum discord is employed now [3,4,5]. The basis for the discord conception is the idea of measurements performed on a system and maximum amount of classical information being extracted with their help.

Due to the fact that it is necessary to solve the optimization problem, the evaluation of quantum correlations, especially discord, is extremely hard [6]. If for the two-qubit systems the quantum entanglement of formation has been obtained for the arbitrary density matrices [7], the analytical formulas for the quantum discord were proposed for X states [8,9,10,11,12,13]. In an X matrix, nonzero entries may belong only to the main diagonal and anti-diagonal [14,15]. Notice that the sum and product of X matrices is again the X matrix (i.e., a set of X matrices is algebraically closed).

However, it was found later that the formulas [9,10,11,12] are incorrect in general. The reason is that the authors [9,10,11,12] believed (and it was their error) that the optimal measurements are achieved only in the limiting points, i.e., at the angles $\theta = 0$ or $\pi/2$ (see below). But on the explicit examples [16,17,18] of X density matrices it was proved that the optimal measurements can take place at the intermediate angles in the interval $(0, \pi/2)$. Unfortunately, these examples with density matrices are specific and do not clarify the general situation.

In the present paper we show that the domain of intermediate optimal angles can arise in the vicinity of transition from the domain with optimal measurement angle $\theta = \pi/2$ to the domain with optimal angle $\theta = 0$ (or inversely). We derive the equations for the boundaries between these domains and investigate their solutions for different models. In particular, the boundaries can coincide or be absent at all and then the quantum discord is given in the total domain of definition by closed analytical formulas.

In the following sections, the general seven-parameters X density matrix is reduced to the five-parameter form by using local unitary transformations, the existence of intermediate subdomains with the optimal angles $\theta \neq 0, \pi/2$ is proved, the equations for boundaries between different subdomains are derived and then applied to various physical systems. Finally, a brief conclusion is given.

2 Real non-negative form for the X density matrices. Domain of definition for the matrix elements

In the most general case, the X density matrix of two-qubit (A and B or 1 and 2) system is given as

$$\rho_{AB} = \begin{pmatrix} a & 0 & 0 & u_1 + iu_2 \\ 0 & b & v_1 + iv_2 & 0 \\ 0 & v_1 - iv_2 & c & 0 \\ u_1 - iu_2 & 0 & 0 & d \end{pmatrix}, \quad (1)$$

where $a + b + c + d = 1$. This matrix contains seven real parameters which, due to the non-negativity definition of any density operator, must satisfy the inequalities

$$a, b, c, d \geq 0, \quad ad \geq u_1^2 + u_2^2, \quad bc \geq v_1^2 + v_2^2. \quad (2)$$

The quantum entanglement and quantum discord are invariant under the local unitary transformations of density matrices [1,2,3,4,5]. Owing to this property, one can with the help of transformation

$$U = e^{-i\varphi_1\sigma_z/2} \otimes e^{-i\varphi_2\sigma_z/2} \quad (3)$$

reduce the seven-parameters density matrix (1) to the real five-parameters X form [18,19,20,21], $U\rho_{AB}U^\dagger \rightarrow \rho_{AB}$. This provides with the angles

$$\varphi_{1,2} = \frac{1}{2}(\arctan \frac{u_2}{u_1} \pm \arctan \frac{v_2}{v_1}). \quad (4)$$

After this, the density matrix (1) takes the form

$$\rho_{AB} = \begin{pmatrix} a & 0 & 0 & u \\ 0 & b & v & 0 \\ 0 & v & c & 0 \\ u & 0 & 0 & d \end{pmatrix}, \quad (5)$$

where

$$u = u_1 \cos(\arctan \frac{u_2}{u_1}) + u_2 \sin(\arctan \frac{u_2}{u_1}), \quad (6)$$

$$v = v_1 \cos(\arctan \frac{v_2}{v_1}) + v_2 \sin(\arctan \frac{v_2}{v_1}). \quad (7)$$

Moreover, with the help of local rotations again around the z axis, it is not difficult to obtain also the non-negative off-diagonal elements of the X matrix (5). Indeed, the local unitary transformation

$$U_1 = e^{-i\frac{\pi}{4}\sigma_z} \otimes e^{i\frac{\pi}{4}\sigma_z} = \begin{pmatrix} i & & & \\ & 1 & & \\ & & 1 & \\ & & & -i \end{pmatrix}, \quad (8)$$

changes the sign of the off-diagonal matrix element u :

$$U_1 \rho_{AB} U_1^+ = \begin{pmatrix} a & 0 & 0 & -u \\ 0 & b & v & 0 \\ 0 & v & c & 0 \\ -u & 0 & 0 & d \end{pmatrix}. \quad (9)$$

Similarly, the local transformation

$$U_2 = e^{i\frac{\pi}{4}\sigma_z} \otimes e^{-i\frac{\pi}{4}\sigma_z} = \begin{pmatrix} 1 & & & \\ & i & & \\ & & -i & \\ & & & 1 \end{pmatrix}, \quad (10)$$

selectively acts on the sign of v :

$$U_2 \rho_{AB} U_2^+ = \begin{pmatrix} a & 0 & 0 & u \\ 0 & b & -v & 0 \\ 0 & -v & c & 0 \\ u & 0 & 0 & d \end{pmatrix}. \quad (11)$$

Thus, after transformation of X matrix to the real form we may simply enclose the off-diagonal elements in the modul symbols:

$$\rho_{AB} = \begin{pmatrix} a & 0 & 0 & |u| \\ 0 & b & |v| & 0 \\ 0 & |v| & c & 0 \\ |u| & 0 & 0 & d \end{pmatrix}. \quad (12)$$

This operation does not influence on the value of quantum correlations in the system.

Thus, one can now consider the density matrices (5) with restrictions

$$a, b, c, d, u, v \geq 0, \quad a + b + c + d = 1, \quad ad \geq u^2, \quad bc \geq v^2. \quad (13)$$

These relations define the domain \mathcal{D} of X density matrix in the space of its entries.

We can rewrite the density matrix (5) in the equivalent form

$$\rho_{AB} = \frac{1}{4} \begin{pmatrix} 1 + s_1 + s_2 + c_3 & . & . & c_1 - c_2 \\ . & 1 + s_1 - s_2 - c_3 & c_1 + c_2 & . \\ . & c_1 + c_2 & 1 - s_1 + s_2 - c_3 & . \\ c_1 - c_2 & . & . & 1 - s_1 - s_2 + c_3 \end{pmatrix}, \quad (14)$$

where

$$\begin{aligned} s_1 &= a + b - c - d, & s_2 &= a - b + c - d, \\ c_1 &= 2(v + u), & c_2 &= 2(v - u), & c_3 &= a - b - c + d. \end{aligned} \quad (15)$$

Decomposition of this matrix on the Pauli matrices σ_α ($\alpha = x, y, z$) leads to its Bloch form

$$\rho_{AB} = \frac{1}{4}(1 + s_1\sigma_z \otimes 1 + s_21 \otimes \sigma_z + c_1\sigma_x \otimes \sigma_x + c_2\sigma_y \otimes \sigma_y + c_3\sigma_z \otimes \sigma_z). \quad (16)$$

The expansion coefficients are the unary and binary correlation functions and therefore five parameters of density matrix are expressed through the five different correlators,

$$\begin{aligned} s_1 &= \langle \sigma_z^1 \rangle = \text{Tr}(\rho_{AB}\sigma_z \otimes 1), & s_2 &= \langle \sigma_z^2 \rangle = \text{Tr}(\rho_{AB}1 \otimes \sigma_z), \\ c_1 &= \langle \sigma_x^1 \sigma_x^2 \rangle = \text{Tr}(\rho_{AB}\sigma_x \otimes \sigma_x), & c_2 &= \langle \sigma_y^1 \sigma_y^2 \rangle = \text{Tr}(\rho_{AB}\sigma_y \otimes \sigma_y), \\ c_3 &= \langle \sigma_z^1 \sigma_z^2 \rangle = \text{Tr}(\rho_{AB}\sigma_z \otimes \sigma_z). \end{aligned} \quad (17)$$

It is clear that

$$-1 \leq s_1, s_2, c_1, c_2, c_3 \leq 1. \quad (18)$$

The domain of definition, \mathcal{D} , in the space $(s_1, s_2, c_1, c_2, c_3)$ is formed, according to Eqs. (13) and (15), by conditions (see also [17,22])

$$(1 - c_3)^2 \geq (s_1 - s_2)^2 + (c_1 + c_2)^2, \quad (1 + c_3)^2 \geq (s_1 + s_2)^2 + (c_1 - c_2)^2. \quad (19)$$

The solid \mathcal{D} is finite and lies in the five-dimensional hypercube (18). Numerical calculations show that the volume of \mathcal{D} is 8% of the hypercube one.

The domain \mathcal{D} is bounded by two quadratic hypersurfaces

$$(s_1 - s_2)^2 + (c_1 + c_2)^2 - (c_3 - 1)^2 = 0 \quad (20)$$

and

$$(s_1 + s_2)^2 + (c_1 - c_2)^2 - (c_3 + 1)^2 = 0. \quad (21)$$

Rotation by the angle $\pi/4$ around the c_3 axis transforms these hyperquadrics to the forms

$$(s'_2)^2 + (c'_1)^2 - \frac{(c_3 - 1)^2}{2} = 0 \quad (22)$$

and

$$(s'_1)^2 + (c'_2)^2 - \frac{(c_3 + 1)^2}{2} = 0, \quad (23)$$

where

$$s'_{1,2} = (\pm s_1 + s_2)/\sqrt{2}, \quad c'_{1,2} = (\pm c_1 + c_2)/\sqrt{2}. \quad (24)$$

Thus, the five-dimensional domain \mathcal{D} results from an intersection of two conic hypercylinders (22) and (23).

3 Three alternatives for the quantum discord

As mentioned above, the measurement operations lie in the ground of discord notion. Following the founders of discord conception [23,24] and their adherents [8,9,10,11,12] we will consider here the one-dimensional projective measurements (i. e., the von Neumann measurements). Such measurements can be reduced to projections which are characterized by the polar (θ) and azimuthal (ϕ) angles relative to the z axis. It is important that in the case of real X density matrix with an additional condition $uv \geq 0$ the optimal measurements are achieved by $\cos 2\phi = 1$ [18,19]. Since the sign of off-diagonal elements are changed by the local unitary transformations, we can always satisfy the above condition.

In the general case, quantum discord depends on subsystem (A or B) where the measurements are performed. For definiteness and without loss of generality, let the measured subsystem be B . Then the quantum discord is given as [3,4,5]

$$Q = S(\rho_B) - S(\rho_{AB}) + \min_{\theta} S_{cond}(\theta), \quad (25)$$

where $\rho_B = \text{Tr}_A \rho_{AB}$ is the reduced density matrix and $S(\rho) = -\text{Tr} \rho \ln \rho$ is the von Neumann entropy for the corresponding state ρ . (Here the entropy is in nats; to transform it, e.g., in bits, one should divide it by $\ln 2$.) Simple calculations with (5) lead to

$$S(\rho_B) = -(a+c) \ln(a+c) - (b+d) \ln(b+d), \quad (26)$$

$S(\rho_{AB}) = S$, where

$$\begin{aligned} S = & -\frac{a+d+\sqrt{(a-d)^2+4u^2}}{2} \ln \frac{a+d+\sqrt{(a-d)^2+4u^2}}{2} \\ & -\frac{a+d-\sqrt{(a-d)^2+4u^2}}{2} \ln \frac{a+d-\sqrt{(a-d)^2+4u^2}}{2} \\ & -\frac{b+c+\sqrt{(b-c)^2+4v^2}}{2} \ln \frac{b+c+\sqrt{(b-c)^2+4v^2}}{2} \\ & -\frac{b+c-\sqrt{(b-c)^2+4v^2}}{2} \ln \frac{b+c-\sqrt{(b-c)^2+4v^2}}{2}. \end{aligned} \quad (27)$$

The quantum conditional entropy of subsystem A is given as [18]

$$S_{cond}(\theta) = A_1 \ln A_1 + A_2 \ln A_2 - \sum_{i=1}^4 \lambda_i \ln \lambda_i, \quad (28)$$

where

$$A_{1,2} = \frac{1}{2} [1 \pm (a-b+c-d) \cos \theta], \quad (29)$$

$$\begin{aligned} \lambda_{1,2} = & \frac{1}{4} [1 + (a-b+c-d) \cos \theta \\ & \pm \{[a+b-c-d + (a-b-c+d) \cos \theta]^2 + 4u^2 \sin^2 \theta\}^{1/2}], \end{aligned} \quad (30)$$

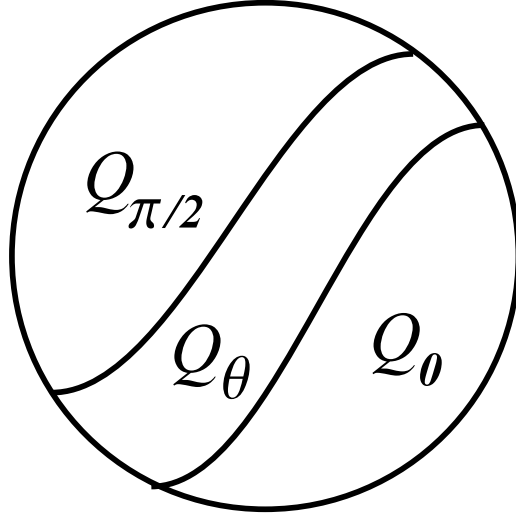


Fig. 1 A fragment of phase diagram with three possible subdomains for the X -state quantum discord

$$\lambda_{3,4} = \frac{1}{4} \llbracket 1 - (a - b + c - d) \cos \theta \pm \{ [a + b - c - d - (a - b - c + d) \cos \theta]^2 + 4w^2 \sin^2 \theta \}^{1/2} \rrbracket, \quad (31)$$

$$w = |u| + |v|. \quad (32)$$

Thus, S_{cond} depends in fact on four parameters because u and v enter via the combination (32). The conditional entropy $S_{cond}(\theta)$ is a differentiable function of its argument θ .

Expressions (26)-(31) allow to define the measurement-dependent discord as [4]

$$Q(\theta) = S(\rho_B) - S(\rho_{AB}) + S_{cond}(\theta), \quad (33)$$

where $\theta \in [0, \pi/2]$. It is obvious that the absolute minimum of this discord can be either on the bounds ($\theta = 0, \pi/2$) or at the intermediate point $\theta \in (0, \pi/2)$. As a result, there is a choice from three possibilities for the quantum discord

$$Q = \min\{Q_0, Q_{\theta}, Q_{\frac{\pi}{2}}\}. \quad (34)$$

This equation generalizes the one proposed earlier for the quantum discord [8, 9, 10, 11, 12]

$$\tilde{Q} = \min\{Q_0, Q_{\frac{\pi}{2}}\}, \quad (35)$$

i.e., it was assumed that the optimal observable can be either σ_z or σ_x . In Fig. 1 we schematically illustrate the parameter domain of a system with three possible subdomains for the discord.

From Eqs. (26)-(33), we have for the discord branch $Q_0 \equiv Q(0)$:

$$Q_0 = -S - a \ln a - b \ln b - c \ln c - d \ln d. \quad (36)$$

For $\theta = \pi/2$ we obtain

$$Q_{\frac{\pi}{2}} = -S - (a+c) \ln(a+c) - (b+d) \ln(b+d) \\ - \frac{1 + \sqrt{(a+b-c-d)^2 + 4w^2}}{2} \ln \frac{1 + \sqrt{(a+b-c-d)^2 + 4w^2}}{2} \\ - \frac{1 - \sqrt{(a+b-c-d)^2 + 4w^2}}{2} \ln \frac{1 - \sqrt{(a+b-c-d)^2 + 4w^2}}{2}. \quad (37)$$

Thus, the branches Q_0 and $Q_{\pi/2}$ are expressed analytically, and the branch $Q_\theta = \min_{\theta \in (0, \pi/2)} Q(\theta)$, if the intermediate minimum exists, should be found from the numerical solution of one-dimensional minimization problem or from the transcendental equation

$$S'_{cond}(\theta) = 0. \quad (38)$$

The first derivative of conditional entropy with respect to θ is equal to

$$S'_{cond}(\theta) = \Lambda'_1(1 + \ln \Lambda_1) + \Lambda'_2(1 + \ln \Lambda_2) - \sum_{i=1}^4 \lambda'_i(1 + \ln \lambda_i) \quad (39)$$

with

$$\Lambda'_{1,2} = \mp \frac{1}{2}(a-b+c-d) \sin \theta, \quad (40)$$

$$\lambda'_{1,2} = \frac{1}{4} \left[-(a-b+c-d) \sin \theta \right. \\ \left. \pm \frac{[a+b-c-d + (a-b-c+d) \cos \theta][-(a-b-c+d) \sin \theta] + 2w^2 \sin 2\theta}{\sqrt{[a+b-c-d + (a-b-c+d) \cos \theta]^2 + 4w^2 \sin^2 \theta}} \right], \quad (41)$$

$$\lambda'_{3,4} = \frac{1}{4} \left[(a-b+c-d) \sin \theta \right. \\ \left. \pm \frac{[a+b-c-d - (a-b-c+d) \cos \theta](a-b-c+d) \sin \theta + 2w^2 \sin 2\theta}{\sqrt{[a+b-c-d - (a-b-c+d) \cos \theta]^2 + 4w^2 \sin^2 \theta}} \right]. \quad (42)$$

All three possible variants for the quantum discord (Q_0 , $Q_{\pi/2}$, and Q_θ) can really exist in physical systems. In the case when $a = b$ and $b = c$ (or $s_1 = s_2 = 0$) the conditional entropy minimum is always achieved at one of two bound points [8]. However, this is wrong for the more general X states; global minimum can take place at inner points of the interval $(0, \pi/2)$.

Indeed, following the authors [16], let us consider the state

$$\rho = \begin{pmatrix} 0.0783 & 0 & 0 & 0 \\ 0 & 0.125 & 0.100 & 0 \\ 0 & 0.100 & 0.125 & 0 \\ 0 & 0 & 0 & 0.6717 \end{pmatrix}. \quad (43)$$

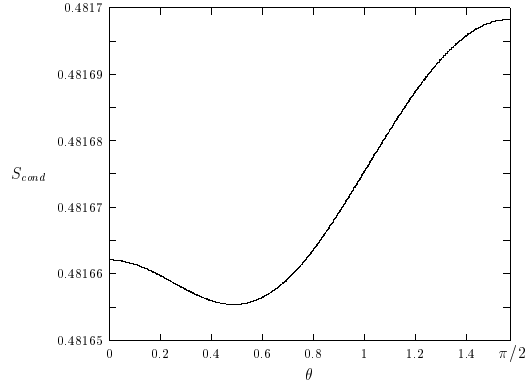


Fig. 2 Quantum conditional entropy S_{cond} as a function of measured angle θ for the state (43)

Using Eqs. (28)-(32) we have computed the function $S_{cond}(\theta)$ for this state. Its behavior is shown in Fig. 2. From the figure we conclude that the conditional entropy minimum is situated in the intermediate region, namely, at the angle $\theta = 0.4883 \approx 28^\circ$. Two other similar numerical examples of quantum states are given in Ref. [18].

These examples clearly show that the optimal measurement angles can really be in the intermediate region $(0, \pi/2)$, i.e., the optimal observables for quantum discord can be not only the σ_x or σ_z , but also their superposition.

For the real X state with constraint $|u + v| \geq |u - v|$ (i.e., $uv \geq 0$ or $\text{sign} u = \text{sign} v$), the authors [17] have proved a theorem which guarantees that the optimal observable is σ_z if

$$(|u| + |v|)^2 \leq (a - b)(d - c) \quad (44)$$

and σ_x if

$$|u| + |v| \geq |\sqrt{ad} - \sqrt{bc}|. \quad (45)$$

The theorem states nothing for the region lying between these bounds. But in the case

$$ac = bd \quad (46)$$

the inequalities (44) and (45) lead to absence of the intermediate region [25].

4 Equations for the boundaries

Let us start with a heuristic example. Consider a two-parameter family of X states [17, 26]

$$\rho = \begin{pmatrix} \epsilon/2 & 0 & 0 & \epsilon/2 \\ 0 & (1-\epsilon)m & 0 & 0 \\ 0 & 0 & (1-\epsilon)(1-m) & 0 \\ \epsilon/2 & 0 & 0 & \epsilon/2 \end{pmatrix} \quad (47)$$

or

$$\begin{aligned}\rho &= \frac{1}{4}[1 + (1 - \epsilon)(2m - 1)(\sigma_1^z - \sigma_2^z) + \epsilon(\sigma_1^x \sigma_2^x - \sigma_1^y \sigma_2^y) + (2\epsilon - 1)\sigma_1^z \sigma_2^z] \\ &= \epsilon|\Phi^+\rangle\langle\Phi^+| + (1 - \epsilon)m|01\rangle\langle 01| + (1 - \epsilon)(1 - m)|10\rangle\langle 10|,\end{aligned}\quad (48)$$

where $|\Phi^+\rangle = (|00\rangle + |11\rangle)/\sqrt{2}$. The given density matrix ρ represents the generalized Horodecki states [22].

Simple calculation yields $Q_0 = \epsilon$ (in bits). Sufficient conditions (44) and (45) for the Q_0 and $Q_{\pi/2}$ subdomains give [17]

$$\epsilon \leq \frac{2m(1 - m)}{1 + 2m(1 - m)} \quad (49)$$

and

$$\epsilon \geq \frac{\sqrt{m(1 - m)}}{1 + \sqrt{m(1 - m)}}, \quad (50)$$

respectively. But in the region

$$\frac{2m(1 - m)}{1 + 2m(1 - m)} < \epsilon < \frac{\sqrt{m(1 - m)}}{1 + \sqrt{m(1 - m)}} \quad (51)$$

the above theorem does not say anything.

Let us now find the lines on the plane (m, ϵ) which are defined by the condition

$$Q_0(m, \epsilon) = Q_{\frac{\pi}{2}}(m, \epsilon). \quad (52)$$

Then we will study the changes of curves $S_{cond}(\theta)$ in the neighborhood to those lines.

Using Eqs. (27), (36), and (37) we have numerically solved the transcendental equation (52). The solution is only one. The results are plotted in Fig. 3 by dotted line.

Consider in detail a particular case. Let the ϵ is held fixed and equal, for example, to $\epsilon = 0.228$ (see Fig. 3). Then the equality $Q_0 = Q_{\pi/2}$ is satisfied at the crossing point $m_{\times} = 0.101\,234$. Study now the behavior of $S_{cond}(\theta)$ when the parameter m varies. Inequalities (49) and (50) guarantee that when $m < 0.096\,545$ the discord equals $Q = Q_{\pi/2}$ and $Q = Q_0$ when $m > 0.180\,107$. If $m = 0.1015$, the minimum of $S_{cond}(\theta)$ is at $\theta = 0$ [see Fig. 4 (a)]. Moreover, the angle $\theta = 0$ is optimal for all larger values of m . When the m decreases the minimum on the curve $S_{cond}(\theta)$ inside the interval between 0 and $\pi/2$ appears. The minimum is clearly seen when $m = 0.1014$ [Fig. 4 (b)]. Near the point $m = 0.101\,234$ the minimum achieves large depth. By further decreasing m , the minimum moves to the bound $\theta = \pi/2$ and then it disappears at all. Optimal measurements undergo to the angle $\theta = \pi/2$.

We argue now that both lower and upper boundaries of the interval where the optimal angles lie between 0 and $\pi/2$ are exact, i.e., the intermediate minimum of $S_{cond}(\theta)$ suddenly appears and suddenly disappears. Above all, we note that the first derivative of function $S_{cond}(\theta)$ at $\theta = 0$ and $\pi/2$ equals

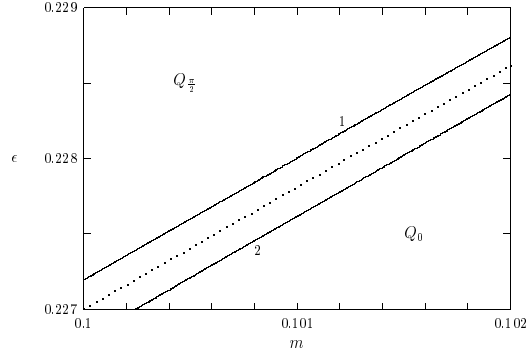


Fig. 3 Subdomains $Q_{\pi/2}$, Q_0 , and (between them) Q_θ for the state (47). Dotted line corresponds to the condition $Q_{\pi/2} = Q_0$. Solid lines 1 and 2 are the $\pi/2$ - and 0-boundaries, respectively

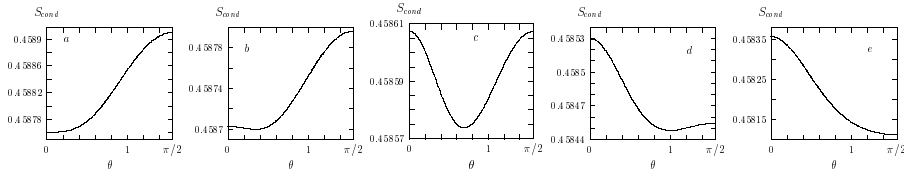


Fig. 4 Appearance and disappearance of an intermediate minimum on the conditional entropy curve by transition from Q_0 to $S_{\pi/2}$ subdomain. Here, $S_{cond}(\theta)$ corresponds to the state (47) at the fixed value of $\epsilon = 0.228$ and $m = 0.1015$ (a), 0.1014 (b), 0.101234 (c), 0.1011 (d), and 0.1008 (e)

zero in general case: $S'_{cond}(0) \equiv S'_{cond}(\pi/2) \equiv 0$. This is easy to check by direct calculations using Eqs. (39)-(42). Let us turn now again to the Fig. 4. By fixed value of parameter ϵ and for each value of m one can say at any moment the inside minimum exists or it is absent. For instance, when $m = 0.1015$ ($\epsilon = 0.228$) the function $S_{cond}(\theta)$ is concave at the point $\theta = 0$ and therefore its second derivative $S''_{cond}(\pi/2) < 0$. But when $m = 0.1014$ the conditional entropy has a local maximum at the same bound point $\theta = 0$ and therefore $S''_{cond}(\pi/2) > 0$. Hence, the bifurcation point (in the sense that two extrema arise from one) [27] is determined by the condition

$$S''_{cond}(0) = 0. \quad (53)$$

Similarly we have for the other bound point $\theta = \pi/2$,

$$S''_{cond}(\pi/2) = 0. \quad (54)$$

Using Eqs. (28)-(31) we get the second derivatives at limiting points:

$$\begin{aligned} S''_{cond}(0) &= \frac{1}{4}(a - b + c - d) \left(2 \ln \frac{b+d}{a+c} + \ln \frac{ac}{bd} \right) \\ &+ \frac{1}{4}(a - b - c + d) \ln \frac{ad}{bc} - \frac{1}{2} w^2 \left(\frac{1}{a-c} \ln \frac{a}{c} + \frac{1}{b-d} \ln \frac{b}{d} \right) \end{aligned} \quad (55)$$

and

$$\begin{aligned}
S''_{cond}(\pi/2) = & \frac{8w^2}{r^3} [(a-c)(b-d) + w^2] \ln \frac{1+r}{1-r} + (a-b+c-d)^2 \\
& - \frac{1}{2(1+r)} \left[a-b+c-d + \frac{1}{r}(a+b-c-d)(a-b-c+d) \right]^2 \\
& - \frac{1}{2(1-r)} \left[a-b+c-d - \frac{1}{r}(a+b-c-d)(a-b-c+d) \right]^2, \quad (56)
\end{aligned}$$

where

$$r = [(a+b-c-d)^2 + 4w^2]^{1/2} \quad (57)$$

and w is given by Eq. (32). The relations (53)-(57) are the boundary equations for the crossover subdomain Q_θ [20,21]. Thus, the boundaries consist of bifurcation points.

If the solutions of Eqs. (53) and (54) are the same, the intermediate subdomain Q_θ is absent and the quantum discord is given by analytical expressions. On the other hand, instead of rough conditions (44) and (45), the inequalities $S''_{cond}(0) \leq 0$ and $S''_{cond}(\pi/2) \leq 0$ define now the complete subdomains Q_0 and $Q_{\pi/2}$, respectively.

Numerical solution of Eqs. (53)-(57) for the state (47) shows that the boundaries are the lines going approximately parallel to the dotted lines (see the lines 1 and 2 in Fig. 3). As a result, the subdomain appears within which the optimal angles should be found numerically. Out of this subdomain we have analytical expressions for the quantum discord. By $\epsilon = 0.228$, the value for m of $\pi/2$ -boundary equals $m_{\pi/2} = 0.100997$ and for the 0-boundary it is $m_0 = 0.101474$. The middle of this interval equals 0.101236 which is near the point $m_\times = 0.101234$.

Consider the discord behavior by a transition from the subdomain $Q_{\pi/2}$ to Q_0 one (Fig. 5). One can see that down to crossing point $m_\times = 0.101234$ the discord \tilde{Q} , according to Refs. [9,10,11], equals $Q_{\pi/2}$ and above the point m_\times it equals Q_0 (see Fig. 5). If this were valid, the discord $\tilde{Q} = \min\{Q_{\pi/2}, Q_0\}$ would not be differentiable at the intersection point m_\times . However, in fact, the true discord $Q = \min\{Q_{\pi/2}, Q_\theta, Q_0\}$ is smooth. This follows from the numerical solution of the task in the intermediate domain. The results are shown again in Fig. 5 by solid line. It is clearly seen that smoothness occurs. We may say that, instead a fracture at m_\times , two hidden transitions occur at the $\pi/2$ - and 0-boundaries.

Notice that the conditions (49) and (50) are rough too and lead to the bounds which lie far beyond the region of Fig. 3.

5 Bell-diagonal states

The case $a = d$ and $b = c$ or $s_1 = s_2 = 0$ (i.e., when both local Bloch vectors are zero) corresponds to the Bell-diagonal states. Domain of definition for the physical states, \mathcal{D} , lies now in the three-dimensional cube defined by $c_1, c_2, c_3 \in$

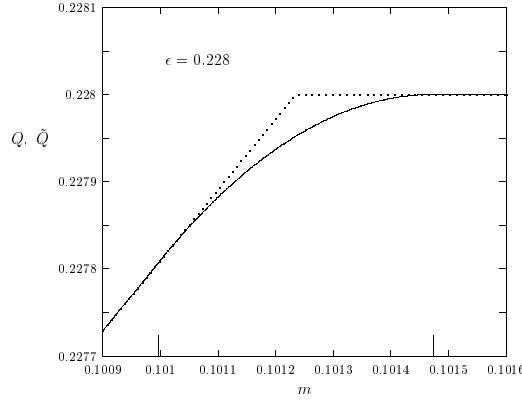


Fig. 5 Dependencies of the false discord $\tilde{Q} = \min\{Q_{\pi/2}, Q_0\}$ (dotted line) and the corrected quantum discord $Q = \min\{Q_{\pi/2}, Q_\theta, Q_0\}$ (solid line) for the state (47) with parameter $\epsilon = 0.228$. Longer bars mark the exact boundaries $m_{\pi/2} = 0.100997$ and $m_0 = 0.101474$. Subdomains $m \leq m_{\pi/2}$, $m_{\pi/2} < m < m_0$, and $m \geq m_0$ correspond to the discord branches $Q_{\pi/2}$, Q_θ , and Q_0 respectively

$[-1, 1]$. Two second-order hypersurfaces (20) and (21) are transformed to the two first-order surfaces

$$\pm |c_1 + c_2| + c_3 - 1 = 0 \quad (58)$$

and

$$\pm |c_1 - c_2| + c_3 + 1 = 0. \quad (59)$$

The former consists of two semi-planes with a \wedge -shaped cross section and the latter is similar to it but has a \vee -shaped cross section. The angle between semi-planes equals $\arccos(1/3) \approx 78^\circ$. These semi-plane surfaces put bounds to the domain \mathcal{D} that is reduced, as shown in Fig. 6, to a tetrahedron with vertices [28]

$$v_1 = (-1, 1, 1), \quad v_2 = (1, -1, 1), \quad v_3 = (1, 1, -1), \quad v_4 = (-1, -1, -1); \quad (60)$$

these vertices lie in octants II $(-, +, +)$, IV $(+, -, +)$, V $(+, +, -)$, and VII $(-, -, -)$, respectively. The centers of tetrahedron facets are

$$\begin{aligned} o_1 &= (1/3, 1/3, 1/3), & o_2 &= (-1/3, -1/3, 1/3), \\ o_3 &= (-1/3, 1/3, -1/3), & o_4 &= (1/3, -1/3, -1/3). \end{aligned} \quad (61)$$

Tetrahedron volume equals a third (i.e., about 33.3%) of the cube one. Notice that the tetrahedron vertices are the states with maximal value of discord (which equals one in bit units).

It is known [29] that the states with zero discord are negligible in the whole Hilbert space. In particular, it has been proved [30, 31] that, when $s_1 = s_2 = 0$, the zero-discord states have at most one nonzero component of vector (c_1, c_2, c_3) , i.e., all classical-only correlated states lie on the Cartesian axes

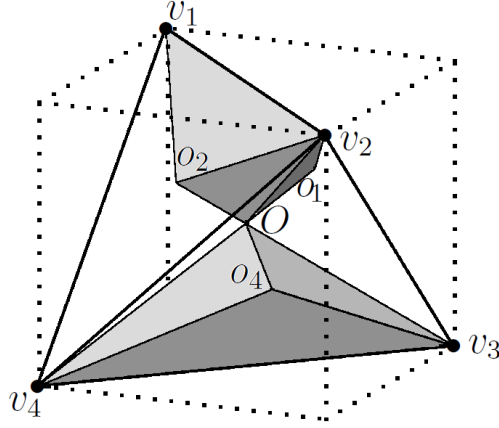


Fig. 6 Tetrahedron with vertices v_1, v_2, v_3 and v_4 is the domain for the total Bell-diagonal states. Two regions (O, v_1, v_2, o_1, o_2) and (O, v_3, v_4, o_3, o_4) correspond to the physical states with Q_0 discord

Oc_1, Oc_2 or Oc_3 . (This corresponds to the so-called “Ising spins” introduced as a matter of fact by his adviser W. Lenz in 1920 [32,33].)

In the case of Bell-diagonal states, both boundary equations (53)-(57) are reduced to a relation

$$(a - b)^2 = (|u| + |v|)^2, \quad (62)$$

so that

$$2|c_3| = |c_1 + c_2| + |c_1 - c_2|. \quad (63)$$

Thus, the $\pi/2$ - and 0-boundaries are coincident, the Q_θ subdomain is absent here, and the quantum discord is given by the explicit analytical formula $Q = \min\{Q_0, Q_{\pi/2}\}$ which is in full agreement with Luo’s results [8].

From Eq. (63), four equations follow

$$c_3 = \pm c_1 \quad c_3 = \pm c_2. \quad (64)$$

These planes divide the tetrahedron into subdomains Q_0 and $Q_{\pi/2}$, where the quantum discord takes the values Q_0 or $Q_{\pi/2}$. Q_0 subdomain consists of two hexahedrons (O, v_1, v_2, o_1, o_2) and (O, v_3, v_4, o_3, o_4) ; they are shown in Fig. 6. The remaining volume of a tetrahedron belongs to the $Q_{\pi/2}$ states. It is in two times larger than the volume of Q_0 states.

The behavior of quantum discord for the Bell-diagonal states along different trajectories is illustrated in Fig. 7 by solid lines. Figure 7 *a* shows the discord as a function of $c_3 \in [-0.95, 0.45]$ by fixed values of $c_1 = 0.3$ and $c_2 = 0.25$. The curve is continuous but has the fractures at $c_3 = \pm 0.3$. They happen when the trajectory crosses the planes dividing the Q_0 and $Q_{\pi/2}$ subdomains (see Fig. 6). In this case, the optimal measurement angle θ varies discontinuously, namely, it jumps from $\theta = 0$ to $\theta = \pi/2$ or inversely. In the vicinity of cross points, the conditional entropy $S_{cond}(\theta)$ changes its form going through a straight line

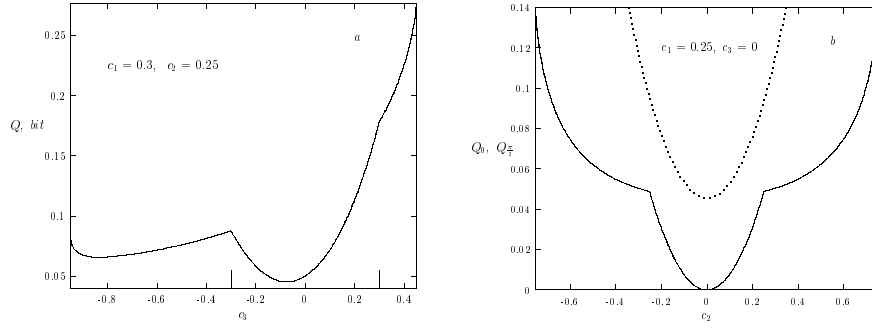


Fig. 7 Quantum discord for the Bell-diagonal states: (a), $Q = \min\{Q_0, Q_{\pi/2}\}$ vs c_3 by $c_1 = 0.3$ and $c_2 = 0.25$, longer bars mark the positions of fracture points at $c_3 = \pm 0.3$; (b), $Q = Q_{\pi/2}$ (solid line) and Q_0 (dotted line) vs c_2 when $c_1 = 0.25$ and $c_3 = 0$

(where any angle $\theta \in [0, \pi/2]$ is optimal). Such a regime of conditional entropy behavior is shown in Fig. 8.

Figure 7 b shows the behavior of branches Q_0 and $Q_{\pi/2}$ as functions of c_2 by fixed values of other two parameters, $c_1 = 0.25$ and $c_3 = 0$. Since here $Q_{\pi/2} < Q_0$, the quantum discord Q equals $Q_{\pi/2}$. The curve $Q_{\pi/2}$ has two fractures. This means that the branch $Q_{\pi/2}$ is a piecewise-analytic function. In this case, however, the optimal measurement angle does not change its value $\theta = \pi/2$ and therefore the position of conditional entropy minimum remains immutable.

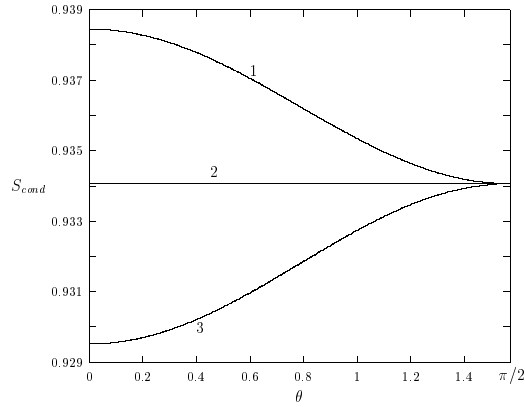


Fig. 8 Transition between $Q_{\pi/2}$ and Q_0 subdomains via a strait line for the conditional entropy. Here, $S_{cond}(\theta)$ is at and near the fracture point $c_3 = 0.3$ on the quantum discord curve in Fig. 7 a. The curves 1, 2, and 3 correspond to $c_3 = 0.29, 0.3$, and 0.31 , respectively

6 Physical systems with the Q_θ subdomains

We are interested now in the systems with Q_θ subdomains. As it was seen from the previous section, such regions do not exist in the Bell-diagonal states. Therefore in this section we will consider systems with nonzero Bloch vectors.

6.1 Phase flip channels

Let us consider the dynamics of quantum discord under decoherence (for a recent review, see, e.g., [34] and references therein). The authors [11] have considered such a dynamics in the phase flip channel. The problem is to calculate the quantum discord for the X matrix

$$\begin{aligned} \varepsilon = \frac{1}{4} [& 1 + s_1 \sigma_z \otimes 1 + s_2 1 \otimes \sigma_z + (1-p)^2 c_1 \sigma_x \otimes \sigma_x \\ & + (1-p)^2 c_2 \sigma_y \otimes \sigma_y + c_3 \sigma_z \otimes \sigma_z]. \end{aligned} \quad (65)$$

Here, the parametrized time $p = 1 - \exp(-\gamma t)$, where t is the time and γ is the phase damping rate. The authors [11] restricted themselves to the case where

$$c_2 = -c_3 c_1, \quad s_2 = c_3 s_1, \quad -1 \leq c_3 \leq 1, \quad -1 \leq s_1 \leq 1. \quad (66)$$

Expansion coefficients in Eq. (65) are related to the corresponding X matrix elements as

$$\begin{aligned} a &= (1 + s_1 + s_2 + c_3)/4, & b &= (1 + s_1 - s_2 - c_3)/4, \\ c &= (1 - s_1 + s_2 - c_3)/4, & d &= (1 - s_1 - s_2 + c_3)/4, \\ u &= (1-p)^2(c_1 - c_2)/4, & v &= (1-p)^2(c_1 + c_2)/4. \end{aligned} \quad (67)$$

Owing to the relation $s_2 = c_3 s_1$, the matrix elements a, b, c , and d satisfy the condition (46) and hence the Q_θ domain is absent here; conditional entropy behaves similar to that as shown in Fig. 8. Thus, nonzero values of s_1 and s_2 are the necessary but not sufficient condition for existence of Q_θ phase.

Consider a different initial state. For example, let us take $s_1 = s_2 = 0.65$, $c_1 = c_2 = 0.249$, and $c_3 = 0.5$. As seen from Fig. 9, the curves $Q_0(p)$ and $Q_{\pi/2}(p)$ have a crossing point at $p_\times \simeq 3158$. An additional study shows that the transition $Q_{\pi/2} \rightarrow Q_0$ goes through the appearance of single minimum on the $S_{cond}(\theta)$ curves inside the interval between 0 and $\pi/2$ (similarly to the curves on Fig. 4).

Solution of equations for the boundaries, Eqs. (53)-(57), shows that the $\pi/2$ - and 0-boundaries do not coincide now and therefore the Q_θ region exists here (see Fig. 10).

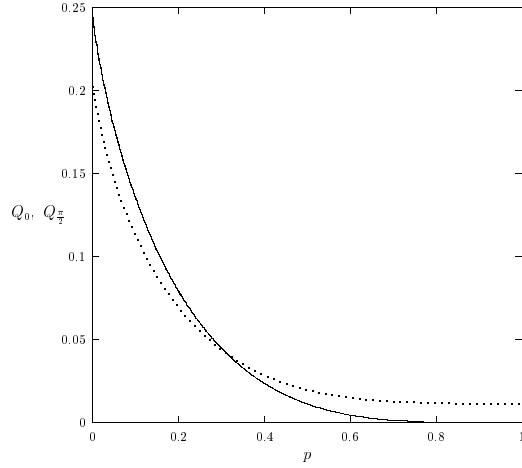


Fig. 9 Q_0 (solid line) and $Q_{\pi/2}$ (dotted line) in bits vs p for the phase flip channel with parameters $s_1 = s_2 = 0.65$, $c_1 = c_2 = 0.249$, and $c_3 = 0.5$. Crossing point of the lines is at $p_{\times} = 0.315789\dots$

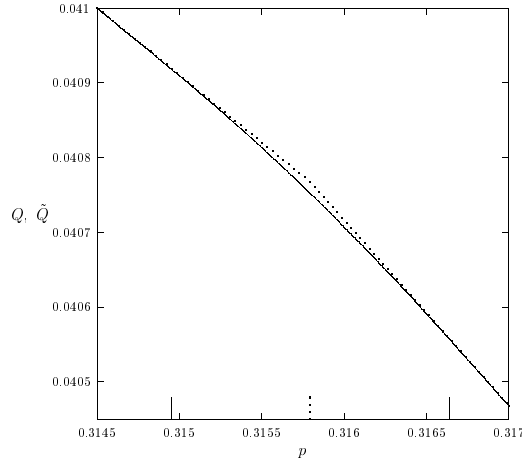


Fig. 10 Dependencies of the false discord $\tilde{Q} = \min\{Q_{\pi/2}, Q_0\}$ (dotted line) and the corrected quantum discord $Q = \min\{Q_{\pi/2}, Q_{\theta}, Q_0\}$ (solid line) vs p for the phase flip channel with parameters $s_1 = s_2 = 0.65$, $c_1 = c_2 = 0.249$, and $c_3 = 0.5$. Longer solid bars mark the boundaries $p_{\pi/2} = 0.314949$ and $p_0 = 0.316637$. Longer dotted bar marks the position of a fracture, $p_{\times} = 0.315789$, on the curve $\tilde{Q}(p)$

6.2 Thermal discord

We now discuss systems at thermal equilibrium. Let us consider the XYZ spin Hamiltonian

$$\mathcal{H} = -\frac{1}{2}(J_x \sigma_1^x \sigma_2^x + J_y \sigma_1^y \sigma_2^y + J_z \sigma_1^z \sigma_2^z + B_1 \sigma_1^z + B_2 \sigma_2^z). \quad (68)$$

This Hamiltonian contains five independent parameters $J_x, J_y, J_z, B_1, B_2 \in (-\infty, \infty)$ (i.e., in \mathcal{R}^5) and is the most general real symmetric traceless X matrix. The corresponding Gibbs density matrix is given as

$$\rho_{AB} = \frac{1}{Z} e^{-\beta \mathcal{H}} \quad (69)$$

(here $\beta = 1/T$, T is the temperature in energy units, Z is the partition function) and has also the five-parameter real X structure. Thus, the map $(B_1/T, B_2/T, J_x/T, J_y/T, J_z/T) \leftrightarrow (s_1, s_2, c_1, c_2, c_3)$ (that is $\mathcal{R}^5 \leftrightarrow \mathcal{D}$) allows in general to change the density-matrix language on a picture of interactions in the XYZ dimer in inhomogeneous fields B_1 and B_2 .

Having solved eigenproblem for the Hamiltonian (68) we then find expressions for the thermal density matrix elements

$$\begin{aligned} a &= \frac{1}{2} \frac{\cosh(\beta R_1/2) + [(B_1 + B_2)/R_1] \sinh(\beta R_1/2)}{\cosh(\beta R_1/2) + \exp(-\beta J_z) \cosh(\beta R_2/2)}, \\ b &= \frac{1}{2} \frac{\cosh(\beta R_2/2) + [(B_1 - B_2)/R_2] \sinh(\beta R_2/2)}{\exp(\beta J_z) \cosh(\beta R_1/2) + \cosh(\beta R_2/2)}, \\ c &= \frac{1}{2} \frac{\cosh(\beta R_2/2) - [(B_1 - B_2)/R_2] \sinh(\beta R_2/2)}{\exp(\beta J_z) \cosh(\beta R_1/2) + \cosh(\beta R_2/2)}, \\ d &= \frac{1}{2} \frac{\cosh(\beta R_1/2) - [(B_1 + B_2)/R_1] \sinh(\beta R_1/2)}{\cosh(\beta R_1/2) + \exp(-\beta J_z) \cosh(\beta R_2/2)}, \\ u &= \frac{1}{2} \frac{[(J_x - J_y)/R_1] \sinh(\beta R_1/2)}{\cosh(\beta R_1/2) + \exp(-\beta J_z) \cosh(\beta R_2/2)}, \\ v &= \frac{1}{2} \frac{[(J_x + J_y)/R_2] \sinh(\beta R_2/2)}{\exp(\beta J_z) \cosh(\beta R_1/2) + \cosh(\beta R_2/2)}, \end{aligned} \quad (70)$$

where

$$R_1 = [(B_1 + B_2)^2 + (J_x - J_y)^2]^{1/2}, \quad R_2 = [(B_1 - B_2)^2 + (J_x + J_y)^2]^{1/2}. \quad (71)$$

For the correlations functions (15), we have respectively

$$\begin{aligned} c_{1,2} &= \frac{2}{Z} \llbracket \pm [(J_x - J_y)/R_1] e^{\beta J_z/2} \sinh(\beta R_1/2) \\ &\quad + [(J_x + J_y)/R_2] e^{-\beta J_z/2} \sinh(\beta R_2/2) \rrbracket, \\ c_3 &= \frac{2}{Z} [e^{\beta J_z/2} \cosh(\beta R_1/2) - e^{-\beta J_z/2} \cosh(\beta R_2/2)], \\ s_{1,2} &= \frac{2}{Z} \llbracket [(B_1 + B_2)/R_1] e^{\beta J_z/2} \sinh(\beta R_1/2) \\ &\quad \pm [(B_1 - B_2)/R_2] e^{-\beta J_z/2} \sinh(\beta R_2/2) \rrbracket, \end{aligned} \quad (72)$$

where the partition function equals

$$Z = 2[e^{\beta J_z/2} \cosh(\beta R_1/2) + e^{-\beta J_z/2} \cosh(\beta R_2/2)] \quad (73)$$

and R_1 and R_2 are given again by Eq. (71).

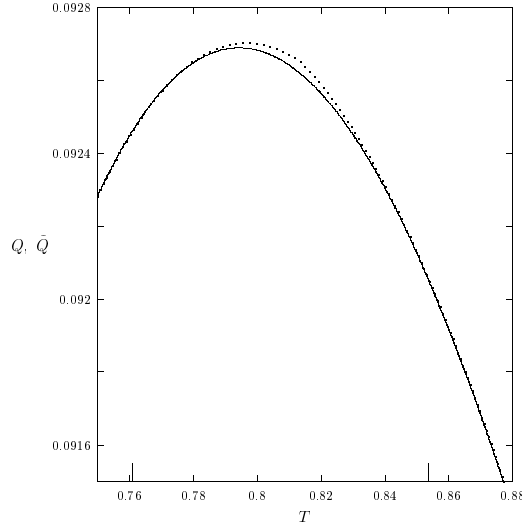


Fig. 11 Dependencies of the false discord $\tilde{Q} = \min\{Q_{\pi/2}, Q_0\}$ (dotted line) and the correct quantum discord $Q = \min\{Q_{\pi/2}, Q_\theta, Q_0\}$ (solid line) for the XXZ dimer with parameters $J = 1$, $J_z = 1.02$ and $B = 1$. Longer bars mark the temperatures $T_{\pi/2} = 0.76106$ and $T_0 = 0.85361$. Domains $T \leq T_{\pi/2}$, $T_{\pi/2} < T < T_0$, and $T \geq T_0$ correspond to the discord branches $Q_{\pi/2}$, Q_θ , and Q_0 , respectively

For each choice of interaction constants J_x , J_y , J_z and external fields B_1 and B_2 we will find the points where the condition $Q_0 = Q_{\pi/2}$ is satisfied. After this we will again study the changes of curves $S_{cond}(\theta)$ in the neighborhood of points found.

Taking, for example, a dimer with parameters $J_x = J_y = J = 1$, $J_z = 1.02$, and $B_1 = B_2 = B = 1$ (that is the XXZ dimer in an uniform field) we consider the thermal discord behavior by a transition from the subdomain $Q_{\pi/2}$ to Q_0 one (Fig. 11). From the figure one can see that down to the crossing point $T_\times = 0.81296$ the discord \tilde{Q} , according to Refs. [9,10,11], equals $Q_{\pi/2}$ and above the point T_\times it equals Q_0 . If this were valid, the discord $\tilde{Q} = \min\{Q_0, Q_{\pi/2}\}$ would have a fracture at the intersection point T_\times . However, in fact, the true discord Q is a smooth function (at least, it is a function of differentiability class C^1). This follows from the numerical solution of the task in the intermediate domain, where the $S_{cond}(\theta)$ curves change similar as in Fig. 4. Results for the quantum discord are shown again in Fig. 11. At the bifurcations points $T_{\pi/2} = 0.76106$ and $T_0 = 0.85361$, the higher derivatives of quantum discord $Q = \min\{Q_{\pi/2}, Q_\theta, Q_0\}$ exhibit a discontinuous behavior.

6.3 Heteronuclear systems with dipolar coupling

Let us consider the system (68) with parameters $J_x = J_y = -D$ and $J_z = 2D$. Such a model corresponds to a dipolar coupled dimer which is stretched along

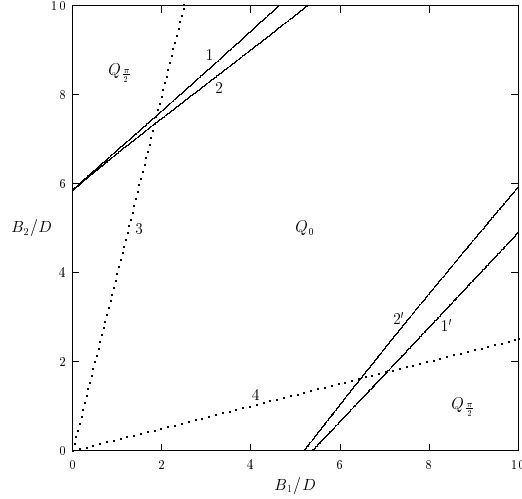


Fig. 12 Subdomains $Q_{\pi/2}$, Q_0 , and (between the lines 1, 2 and 1', 2') Q_{θ} for the spin dimer (74) at the normalized temperature $T/D = 1$. Dotted lines 3 and 4 correspond to $B_2 = 4B_1$ and $B_2 = B_1/4$, respectively

the z axis [35]

$$\mathcal{H} = \frac{1}{2}D(\sigma_1^x \sigma_2^x + \sigma_1^y \sigma_2^y - 2\sigma_1^z \sigma_2^z) - \frac{1}{2}(B_1 \sigma_1^z + B_2 \sigma_2^z). \quad (74)$$

Here the dipolar coupling constant (in frequency units) equals

$$D = \frac{\mu_0}{4\pi} \frac{\gamma_1 \gamma_2}{2r_0^3}, \quad (75)$$

where μ_0 is the magnetic permeability of free space, γ_1 and γ_2 are the gyromagnetic ratios of particles in the dimer, and r_0 is the distance between those particles. Normalized fields B_1 and B_2 in Eq. (75) are

$$B_1 = \gamma_1 B_0, \quad B_2 = \gamma_2 B_0, \quad (76)$$

where B_0 is the external magnetic field induction.

We have performed necessary calculations (according to our approach developed in the previous sections) and found the subdomains of quantum discord in the plane $(B_1/D, B_2/D)$. The results are shown at the normalized temperature $T/D = 1$ in Fig. 12. From this figure one can see that such a system has the Q_{θ} regions (between the 1, 2 and 1', 2' lines). These regions can be reached by varying the external magnetic field B_0 . Two possible trajectories are shown in Fig. 12 by dotted lines, $B_2 = 4B_1$ and $B_2 = B_1/4$. (The value $\gamma_2/\gamma_1 = 4$ approximately corresponds to the quotient of gyromagnetic ratios for the nucleus of ^1H and ^{13}C .)

We found also that in the Q_θ subdomain the conditional entropy $S_{cond}(\theta)$ has only one minimum that is located in the interval $(0, \pi/2)$. The picture is qualitatively similar to that is shown in Fig. 4.

So, the Q_θ region and corresponding sudden changes of quantum correlation behavior at their boundaries can be observed in solid materials with nuclear dimers.

7 Results and perspectives

In light of the above, the calculation of quantum discord of any X states can be achieved by following steps. First, the density matrix (1) is transformed to the real form, i.e., the quantities u and v are calculated using Eqs. (6) and (7). It is also well to solve the equation $Q_0 = Q_{\pi/2}$ and determine possible crossing points of branches Q_0 and $Q_{\pi/2}$. Then the equations $S''_{cond}(0) = 0$ and $S''_{cond}(\pi/2) = 0$ are solved to find the boundaries for the intermediate subdomain Q_θ . After this, one should numerically find the optimal measurement angle $\theta \in (0, \pi/2)$ and compute $Q_\theta = Q(\theta)$. As a result, the quantum discord is given by $Q = \min\{Q_0, Q_\theta, Q_{\pi/2}\}$.

The formula for calculating the quantum discord belongs to a piecewise-defined type

$$f(x) = \begin{cases} F(x, a), & x \in \Omega_a \\ F(x, b), & x \in \Omega_b \\ \min_{\alpha \in (a, b)} F(x, \alpha), & x \in \Omega_c. \end{cases} \quad (77)$$

In other words, the domain of definition, Ω , of the function $f(x)$ consists of subdomains in which the function is given by closed analytical expressions or it exists in a numerical form.

So, the quantum discord of X states is represented analytically if the Q_θ subdomain is absent. Then the quantum discord is given by the closed form $Q = \min\{Q_{\pi/2}, Q_0\}$. The discord is continuous, but generally speaking it is a piecewise smooth function. In particular, this is valid for a spacial class of X states, namely, for the Bell-diagonal states. For them, we found the Q_0 and $Q_{\pi/2}$ regions in the total domain of their definition (Fig. 6). It would be interesting to find the subdomains Q_0 , Q_θ , and $Q_{\pi/2}$ in the five-dimensional domain \mathcal{D} making, e.g., an atlas of maps.

Also, we have shown in this paper that the boundaries for the transition subdomain from Q_0 to $Q_{\pi/2}$ or reversely are exactly defined. They consist of nonanalyticity points which are bifurcation ones. The corresponding equations for these boundaries have been found. The boundaries may coincide and then the quantum discord is evaluated analytically in the total domain of definition. The regions Q_θ with the optimal intermediate angles $\theta \in (0, \pi/2)$ have been found for a number of physical systems including the phase flip channels, spin dimers at the thermal equilibrium, heteronuclear systems with dipolar interaction. The transitions $Q_{\pi/2} \leftrightarrow Q_\theta \leftrightarrow Q_0$ occur continuously and smoothly. This is a new type of transitions for the quantum discord.

We have found only two regimes for the conditional entropy change by above transitions: (i) via the birth of one intermediate minimum (as shown in Fig. 4) and (ii) via the strait line (as shown in Fig. 8). It is hoped that our observations will be rigorously proofed and, maybe, generalized in the future.

At present the attempts are made to obtain analytical formulas for the super quantum discord of X states with nonzero Bloch vectors [36,37]. In this connection one should note the the authors do not take into account a possibility of intermediate optimal angles for the weak measurements which are a generalization of projective ones.

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